Connecting via Winsock to STN

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LOGINID:ssptaeal1624

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * * * * * * Welcome to STN International * * * * * * * * * *

NEWS	1		Web Page for STN Seminar Schedule - N. America
NEWS	2	NOV 21	CAS patent coverage to include exemplified prophetic substances identified in English-, French-, German-, and Japanese-language basic patents from 2004-present
NEWS	3	NOV 26	MARPAT enhanced with FSORT command
NEWS	4	NOV 26	CHEMSAFE now available on STN Easy
NEWS	5	NOV 26	Two new SET commands increase convenience of STN searching
NEWS	6	DEC 01	ChemPort single article sales feature unavailable
NEWS	7	DEC 12	GBFULL now offers single source for full-text coverage of complete UK patent families
NEWS	8	DEC 17	Fifty-one pharmaceutical ingredients added to PS
NEWS	9	JAN 06	The retention policy for unread STNmail messages will change in 2009 for STN-Columbus and STN-Tokyo
NEWS	10	JAN 07	WPIDS, WPINDEX, and WPIX enhanced Japanese Patent Classification Data
NEWS	11	FEB 02	Simultaneous left and right truncation (SLART) added for CERAB, COMPUAB, ELCOM, and SOLIDSTATE
NEWS	12	FEB 02	GENBANK enhanced with SET PLURALS and SET SPELLING
NEWS	13	FEB 06	Patent sequence location (PSL) data added to USGENE
NEWS	14	FEB 10	COMPENDEX reloaded and enhanced

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

NEWS HOURS	STN Operating Hours Plus Help Desk Availability
NEWS LOGIN	Welcome Banner and News Items
NEWS IPC8	For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

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* * * * * * * * * * * * * * * * STN Columbus * * * * * * * * * * * * * * * *

FILE 'HOME' ENTERED AT 16:28:03 ON 10 FEB 2009

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=> file reg
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                         ENTRY    SESSION
FULL ESTIMATED COST                                       0.88       0.88

FILE 'REGISTRY' ENTERED AT 16:30:20 ON 10 FEB 2009
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STRUCTURE FILE UPDATES:    8 FEB 2009  HIGHEST RN 1102960-71-3
DICTIONARY FILE UPDATES:   8 FEB 2009  HIGHEST RN 1102960-71-3

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

  Please note that search-term pricing does apply when
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REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document.  For information
on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

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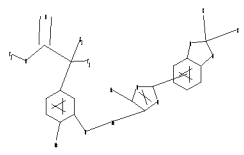

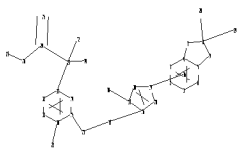

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ring nodes :
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24-26 26-27
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-9 7-8 8-9 10-11 10-15 11-12 12-13 13-14
14-15 16-17 16-20 17-18 18-19 19-20
exact/norm bonds :
10-22 15-21 16-20 16-37 17-30 19-20 21-37 23-34 23-32 24-25 24-26 26-27
exact bonds :
4-7 5-9 7-8 8-9 8-28 8-29 13-23 16-17 17-18 18-19 23-24
normalized bonds :
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isolated ring systems :
containing 1 : 10 : 16 :

G1:C,H

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
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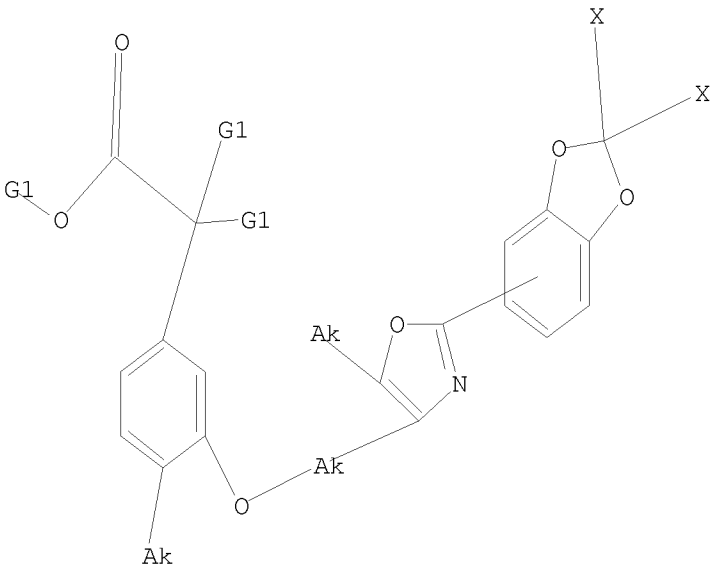

G1 C,H

Structure attributes must be viewed using STN Express query preparation.

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100.0% PROCESSED      130 ITERATIONS                                 4 ANSWERS
SEARCH TIME: 00.00.01

L2            4 SEA SSS FUL L1

=> file caplus
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                                                         ENTRY    SESSION
FULL ESTIMATED COST                                     185.88     186.76

FILE 'CAPLUS' ENTERED AT 16:30:59 ON 10 FEB 2009
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FILE COVERS 1907 - 10 Feb 2009  VOL 150 ISS 7
FILE LAST UPDATED: 9 Feb 2009  (20090209/ED)

Caplus now includes complete International Patent Classification (IPC)
reclassification data for the third quarter of 2008.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate
substance identification.

=> s l2 full
L3            1 L2

=> d ibib abs hitstr tot
```

L3 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2005:283476 CAPLUS
DOCUMENT NUMBER: 142:355258
TITLE: Preparation of azole compounds containing phenylacetic acid moiety as PPARδ agonists
INVENTOR(S): Kusuda, Shinya; Nakayama, Yoshisuke; Tajima, Hisao; Sakamoto, Takahiko
PATENT ASSIGNEE(S): Ono Pharmaceutical Co., Ltd., Japan
SOURCE: PCT Int. Appl., 81 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.		DATE
WO 2005028453	A1	20050331	WO 2004-JP14137		20040921
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RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG					
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IN 2006CN00975	A	20070615	IN 2006-CN975		20060321
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PRIORITY APPLN. INFO.:			JP 2003-330616	A	20030922
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			WO 2004-JP14137	W	20040921

OTHER SOURCE(S): MARPAT 142:355258
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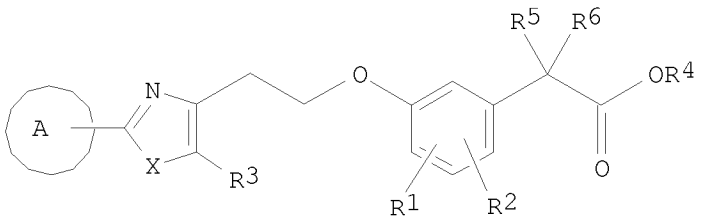

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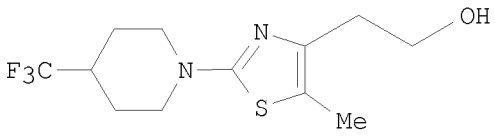

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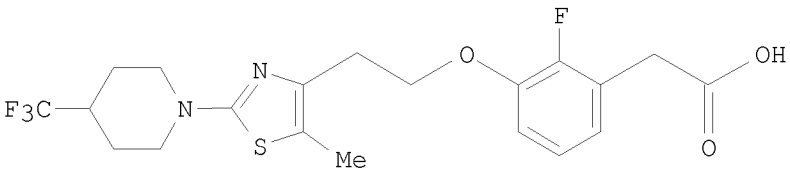


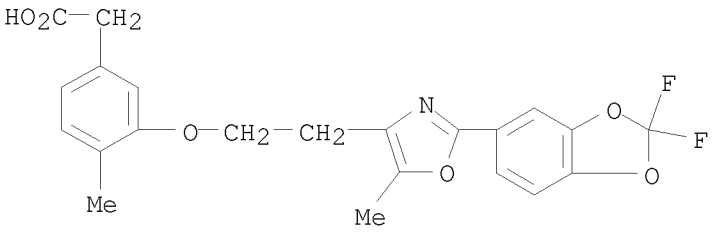
III

AB Title compds. I [R1, R2 = H, alkyl, etc.; R3 = optionally substituted alkyl with halo, etc.; R4 = H, alkyl; R5, R6 = H, alkyl; further detail on R5, R6 is provided.; X = S, O, etc.; ring A = optionally substituted cyclic group] were prepared For example, reaction of compound II, e.g., prepared from 4-(trifluoromethyl)piperidine·HCl in 5 steps, with 2-fluoro-3-hydroxyphenylacetic acid Me ester under Mitsunobu condition followed by hydrolysis using aqueous NaOH afforded compound III. The exemplified compound III exhibited 1.23 fold increase for PPARδ at 1.0 μM. Compds. I are claimed useful as PPARδ agonists for the treatment of hyperlipidemia, obesity. Formulations are given.

IT 848943-42-0P 848943-44-2P 848943-46-4P
848943-47-5P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of azole compds. containing phenylacetic acid moiety as PPAR agonists for treatment of hyperlipidemia, obesity)

RN 848943-42-0 CAPLUS

CN Benzeneacetic acid, 3-[2-[2-(2,2-difluoro-1,3-benzodioxol-5-yl)-5-methyl-4-oxazolyl]ethoxy]-4-methyl- (CA INDEX NAME)



RN 848943-44-2 CAPLUS
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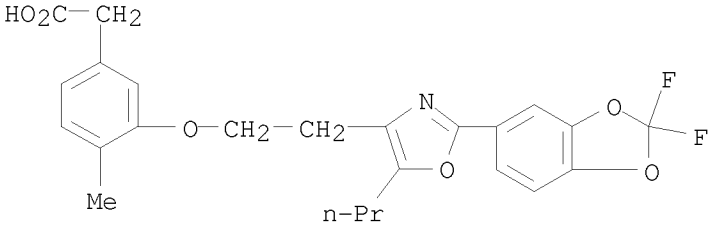

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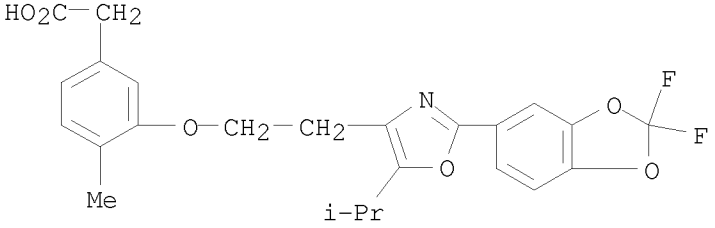

RN 848943-47-5 CAPLUS
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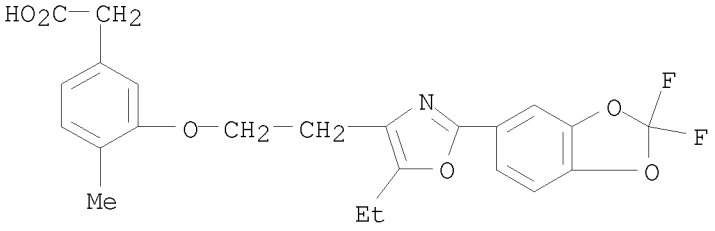

REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS

RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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=> file reg
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
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FULL ESTIMATED COST                                       8.14     194.90

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)          SINCE FILE      TOTAL
                                                         ENTRY    SESSION
CA SUBSCRIBER PRICE                                      -0.82      -0.82

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STRUCTURE FILE UPDATES:    8 FEB 2009  HIGHEST RN 1102960-71-3
DICTIONARY FILE UPDATES:   8 FEB 2009  HIGHEST RN 1102960-71-3

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REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document.  For information
on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

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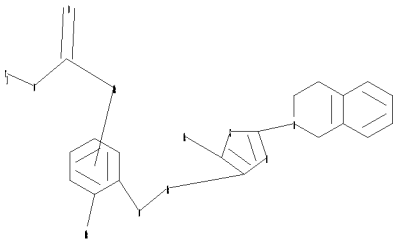

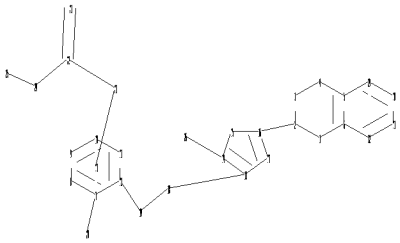

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ring nodes :
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ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-29 6-32 7-8 7-12 8-9 9-10 10-11 11-12 13-14
13-17 14-15 15-16 16-17 29-30 30-31 31-32
exact/norm bonds :
1-2 1-6 2-3 2-16 3-4 4-5 7-20 12-19 13-17 13-18 14-26 16-17 18-19
21-22 22-23 22-24 24-25
exact bonds :
13-14 14-15 15-16
normalized bonds :
5-6 5-29 6-32 7-8 7-12 8-9 9-10 10-11 11-12 29-30 30-31 31-32
isolated ring systems :
containing 1 : 7 : 13 :

G1:C,H

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:CLASS 19:CLASS
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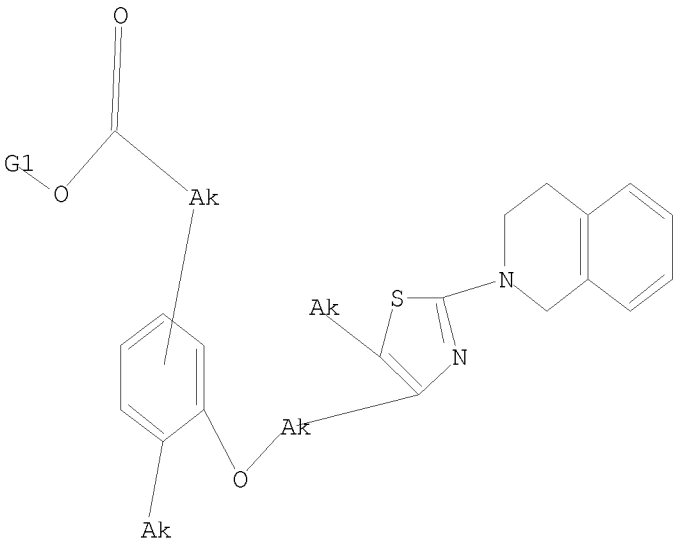

G1 C,H

Structure attributes must be viewed using STN Express query preparation.

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100.0% PROCESSED 53382 ITERATIONS 1 ANSWERS
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L5 1 SEA SSS FUL L4

=> file caplus

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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.82

FILE 'CAPLUS' ENTERED AT 16:34:45 ON 10 FEB 2009

FILE COVERS 1907 - 10 Feb 2009 VOL 150 ISS 7
FILE LAST UPDATED: 9 Feb 2009 (20090209/ED)

Caplus now includes complete International Patent Classification (IPC)
reclassification data for the third quarter of 2008.

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http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate
substance identification.

=> s l5 full
L6 1 L5

=> d ibib abs hitstr

L6 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2005:283476 CAPLUS
DOCUMENT NUMBER: 142:355258
TITLE: Preparation of azole compounds containing phenylacetic acid moiety as PPARδ agonists
INVENTOR(S): Kusuda, Shinya; Nakayama, Yoshisuke; Tajima, Hisao; Sakamoto, Takahiko
PATENT ASSIGNEE(S): Ono Pharmaceutical Co., Ltd., Japan
SOURCE: PCT Int. Appl., 81 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.		DATE
WO 2005028453	A1	20050331	WO 2004-JP14137		20040921
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RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG					
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CA 2539554	A1	20050331	CA 2004-2539554		20040921
EP 1666472	A1	20060607	EP 2004-773449		20040921
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			WO 2004-JP14137	W	20040921

OTHER SOURCE(S): MARPAT 142:355258
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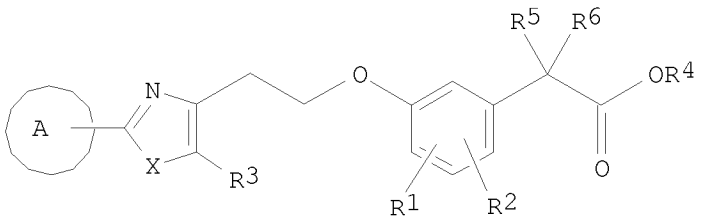

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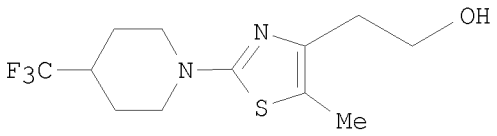

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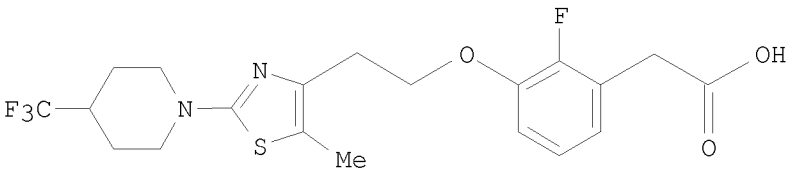


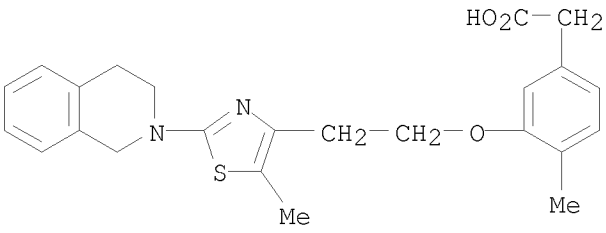
III

AB Title compds. I [R1, R2 = H, alkyl, etc.; R3 = optionally substituted alkyl with halo, etc.; R4 = H, alkyl; R5, R6 = H, alkyl; further detail on R5, R6 is provided.; X = S, O, etc.; ring A = optionally substituted cyclic group] were prepared For example, reaction of compound II, e.g., prepared from 4-(trifluoromethyl)piperidine·HCl in 5 steps, with 2-fluoro-3-hydroxyphenylacetic acid Me ester under Mitsunobu condition followed by hydrolysis using aqueous NaOH afforded compound III. The exemplified compound III exhibited 1.23 fold increase for PPARδ at 1.0 μM. Compds. I are claimed useful as PPARδ agonists for the treatment of hyperlipidemia, obesity. Formulations are given.

IT 848943-61-3P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of azole compds. containing phenylacetic acid moiety as PPAR agonists for treatment of hyperlipidemia, obesity)

RN 848943-61-3 CAPLUS

CN Benzeneacetic acid, 3-[2-[2-(3,4-dihydro-2(1H)-isoquinolinyl)-5-methyl-4-thiazolyl]ethoxy]-4-methyl- (CA INDEX NAME)



REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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FULL ESTIMATED COST                                       8.14     388.92

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CA SUBSCRIBER PRICE                                      -0.82      -1.64

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STRUCTURE FILE UPDATES:    8 FEB 2009  HIGHEST RN 1102960-71-3
DICTIONARY FILE UPDATES:   8 FEB 2009  HIGHEST RN 1102960-71-3

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predicted properties as well as tags indicating availability of
experimental property data in the original document.  For information
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http://www.cas.org/support/stngen/stndoc/properties.html

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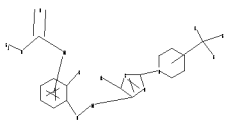

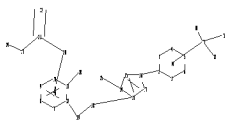

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ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17
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ring bonds :
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exact/norm bonds :
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exact bonds :
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isolated ring systems :
containing 1 : 7 : 13 :

G1:C,H

Match level :
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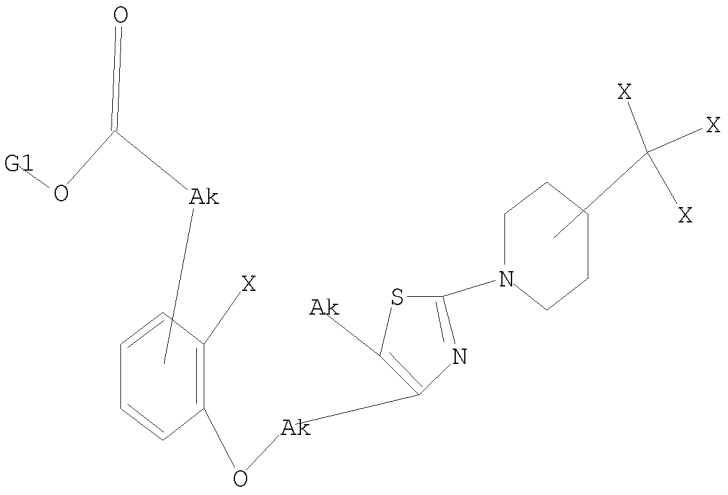

G1 C,H

Structure attributes must be viewed using STN Express query preparation.

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FULL SCREEN SEARCH COMPLETED - 2733 TO ITERATE

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SEARCH TIME: 00.00.01

L8 2 SEA SSS FUL L7

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COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	185.88	574.80

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-1.64

FILE 'CAPLUS' ENTERED AT 16:38:14 ON 10 FEB 2009
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FILE COVERS 1907 - 10 Feb 2009 VOL 150 ISS 7
FILE LAST UPDATED: 9 Feb 2009 (20090209/ED)

Caplus now includes complete International Patent Classification (IPC)
reclassification data for the third quarter of 2008.

This file contains CAS Registry Numbers for easy and accurate
substance identification.

=> s l8 full
L9   2 L8

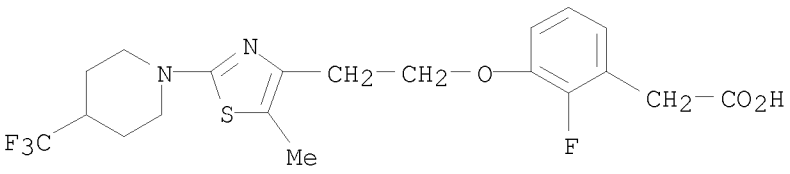
=> d ibib abs hitstr tot

L9 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2007:1388077 CAPLUS
DOCUMENT NUMBER: 149:430
TITLE: Pharmacophore modeling and parallel screening for PPAR ligands
AUTHOR(S): Markt, Patrick; Schuster, Daniela; Kirchmair, Johannes; Laggner, Christian; Langer, Thierry
CORPORATE SOURCE: Department of Pharmaceutical Chemistry, Institute of Pharmacy and Center for Molecular Biosciences Innsbruck (CMBI), University of Innsbruck, Innsbruck, 6020, Austria
SOURCE: Journal of Computer-Aided Molecular Design (2007), 21(10-11), 575-590
CODEN: JCADEQ; ISSN: 0920-654X
PUBLISHER: Springer
DOCUMENT TYPE: Journal
LANGUAGE: English

AB We describe the generation and validation of pharmacophore models for PPARs, as well as a large scale validation of the parallel screening approach by screening PPAR ligands against a large database of structure-based models. A large test set of 357 PPAR ligands was screened against 48 PPAR models to determine the best models for agonists of PPAR-α, PPAR-δ, and PPAR-γ. Afterwards, a parallel screen was performed using the 357 PPAR ligands and 47 structure-based models for PPARs, which were integrated into a 1537 models comprising inhouse pharmacophore database, to assess the enrichment of PPAR ligands within the PPAR hypotheses. For these purposes, we categorized the 1537 database models into 181 protein targets and developed a score that ranks the retrieved targets for each ligand. Thus, we tried to find out if the concept of parallel screening is able to predict the correct pharmacol. target for a set of compds. The PPAR target was ranked first more often than any other target. This confirms the ability of parallel screening to forecast the pharmacol. active target for a set of compds.

IT 848943-49-7
RL: PAC (Pharmacological activity); PRP (Properties); THU (Therapeutic use); BIOL (Biological study); USES (Uses)
(pharmacophore modeling and parallel screening for PPAR ligands)

RN 848943-49-7 CAPLUS
CN Benzeneacetic acid, 2-fluoro-3-[2-[5-methyl-2-[4-(trifluoromethyl)-1-piperidinyl]-4-thiazolyl]ethoxy]- (CA INDEX NAME)



REFERENCE COUNT: 46 THERE ARE 46 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L9 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2005:283476 CAPLUS
DOCUMENT NUMBER: 142:355258
TITLE: Preparation of azole compounds containing phenylacetic acid moiety as PPARδ agonists
INVENTOR(S): Kusuda, Shinya; Nakayama, Yoshisuke; Tajima, Hisao; Sakamoto, Takahiko
PATENT ASSIGNEE(S): Ono Pharmaceutical Co., Ltd., Japan
SOURCE: PCT Int. Appl., 81 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.		DATE
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CN 1882553	A	20061220	CN 2004-80033842		20040921
NO 2006001281	A	20060622	NO 2006-1281		20060321
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MX 2006003205	A	20060623	MX 2006-3205		20060322
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US 20070105868	A1	20070510	US 2006-572937		20060322
PRIORITY APPLN. INFO.:			JP 2003-330616	A	20030922
			JP 2004-231546	A	20040806
			WO 2004-JP14137	W	20040921

OTHER SOURCE(S): MARPAT 142:355258
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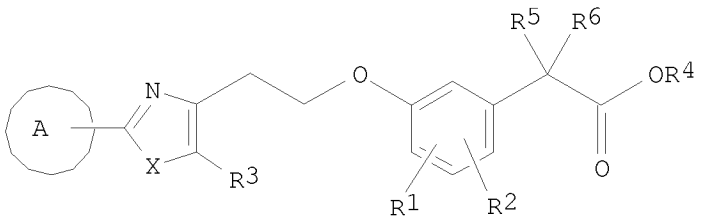

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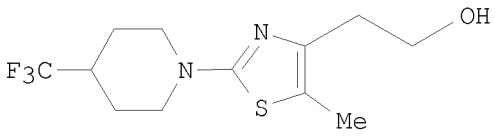

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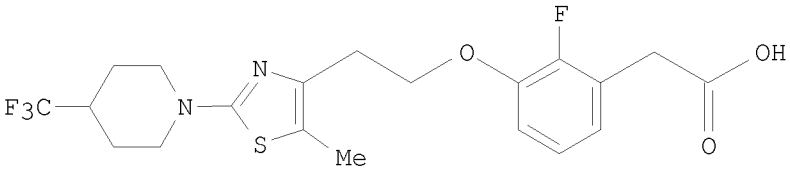


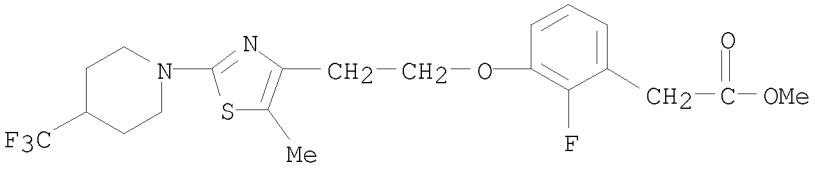
III

AB Title compds. I [R1, R2 = H, alkyl, etc.; R3 = optionally substituted alkyl with halo, etc.; R4 = H, alkyl; R5, R6 = H, alkyl; further detail on R5, R6 is provided.; X = S, O, etc.; ring A = optionally substituted cyclic group] were prepared For example, reaction of compound II, e.g., prepared from 4-(trifluoromethyl)piperidine·HCl in 5 steps, with 2-fluoro-3-hydroxyphenylacetic acid Me ester under Mitsunobu condition followed by hydrolysis using aqueous NaOH afforded compound III. The exemplified compound III exhibited 1.23 fold increase for PPARδ at 1.0 μM. Compds. I are claimed useful as PPARδ agonists for the treatment of hyperlipidemia, obesity. Formulations are given.

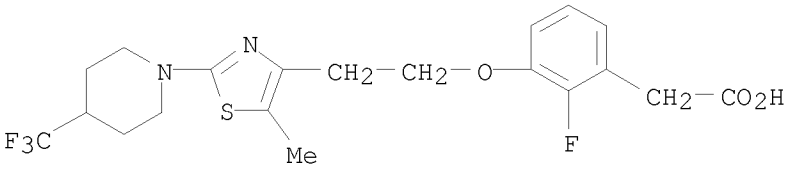
IT 848943-48-6P
RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(preparation of azole compds. containing phenylacetic acid moiety as PPAR agonists for treatment of hyperlipidemia, obesity)

RN 848943-48-6 CAPLUS

CN Benzeneacetic acid, 2-fluoro-3-[2-[5-methyl-2-[4-(trifluoromethyl)-1-piperidinyl]-4-thiazolyl]ethoxy]-, methyl ester (CA INDEX NAME)



IT 848943-49-7P
RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of azole compds. containing phenylacetic acid moiety as PPAR agonists for treatment of hyperlipidemia, obesity)
RN 848943-49-7 CAPLUS
CN Benzeneacetic acid, 2-fluoro-3-[2-[5-methyl-2-[4-(trifluoromethyl)-1-piperidinyl]-4-thiazolyl]ethoxy]- (CA INDEX NAME)



REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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L7             STRUCTURE UPLOADED
L8           2 S L7 FUL

     FILE 'CAPLUS' ENTERED AT 16:38:14 ON 10 FEB 2009
L9           2 S L8 FULL

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'YU' IS NOT VALID HERE
For an explanation, enter "HELP LOGOFF".

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FULL ESTIMATED COST                                  11.78     586.58

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)      SINCE FILE      TOTAL
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STN INTERNATIONAL LOGOFF AT 16:39:01 ON 10 FEB 2009
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